



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:37 AM UTC

PDB ID : 4B3I / pdb_00004b3i
Title : Crystal structure of Mycobacterium tuberculosis fatty acid beta- oxidation complex with CoenzymeA bound at the hydratase active sites
Authors : Venkatesan, R.; Wierenga, R.K.
Deposited on : 2012-07-24
Resolution : 2.63 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

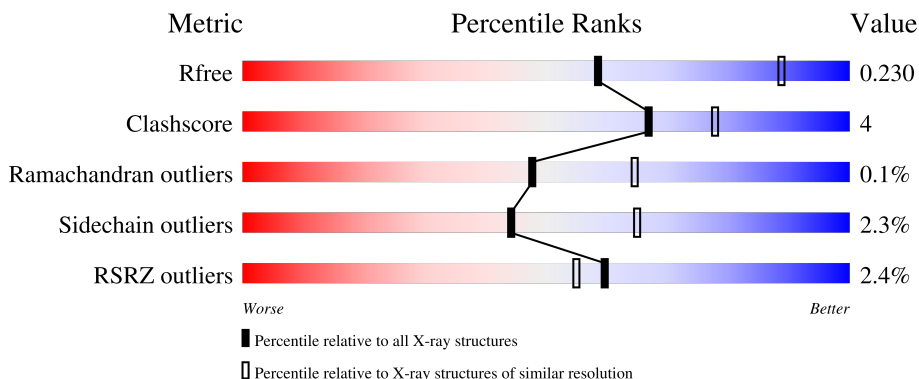
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.63 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	736	% 90% 9% ..
1	B	736	4% 91% 7% .
2	C	403	2% 91% 8%
2	D	403	% 91% 8% .

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17910 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called FATTY ACID BETA-OXIDATION COMPLEX ALPHA-CHAIN FADB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	731	Total	C	N	O	S	0	5	0
			5454	3453	940	1040	21			
1	B	728	Total	C	N	O	S	0	6	0
			5405	3423	925	1037	20			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	expression tag	UNP O53872
A	-14	GLY	-	expression tag	UNP O53872
A	-13	SER	-	expression tag	UNP O53872
A	-12	SER	-	expression tag	UNP O53872
A	-11	HIS	-	expression tag	UNP O53872
A	-10	HIS	-	expression tag	UNP O53872
A	-9	HIS	-	expression tag	UNP O53872
A	-8	HIS	-	expression tag	UNP O53872
A	-7	HIS	-	expression tag	UNP O53872
A	-6	HIS	-	expression tag	UNP O53872
A	-5	SER	-	expression tag	UNP O53872
A	-4	GLN	-	expression tag	UNP O53872
A	-3	ASP	-	expression tag	UNP O53872
A	-2	PRO	-	expression tag	UNP O53872
A	-1	ASN	-	expression tag	UNP O53872
A	0	SER	-	expression tag	UNP O53872
B	-15	MET	-	expression tag	UNP O53872
B	-14	GLY	-	expression tag	UNP O53872
B	-13	SER	-	expression tag	UNP O53872
B	-12	SER	-	expression tag	UNP O53872
B	-11	HIS	-	expression tag	UNP O53872
B	-10	HIS	-	expression tag	UNP O53872
B	-9	HIS	-	expression tag	UNP O53872
B	-8	HIS	-	expression tag	UNP O53872

Continued on next page...

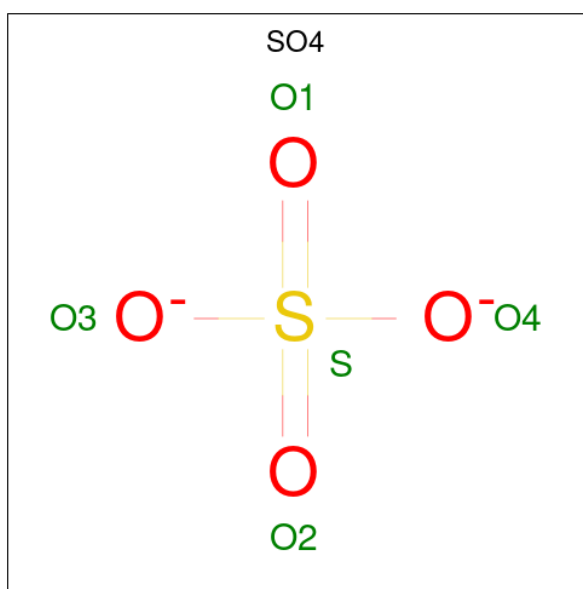
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP O53872
B	-6	HIS	-	expression tag	UNP O53872
B	-5	SER	-	expression tag	UNP O53872
B	-4	GLN	-	expression tag	UNP O53872
B	-3	ASP	-	expression tag	UNP O53872
B	-2	PRO	-	expression tag	UNP O53872
B	-1	ASN	-	expression tag	UNP O53872
B	0	SER	-	expression tag	UNP O53872

- Molecule 2 is a protein called FATTY ACID BETA-OXIDATION COMPLEX BETA-CHAIN FADA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	402	Total	C	N	O	S	0	3	0
			2976	1859	529	573	15			
2	D	403	Total	C	N	O	S	0	2	0
			2968	1854	525	574	15			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

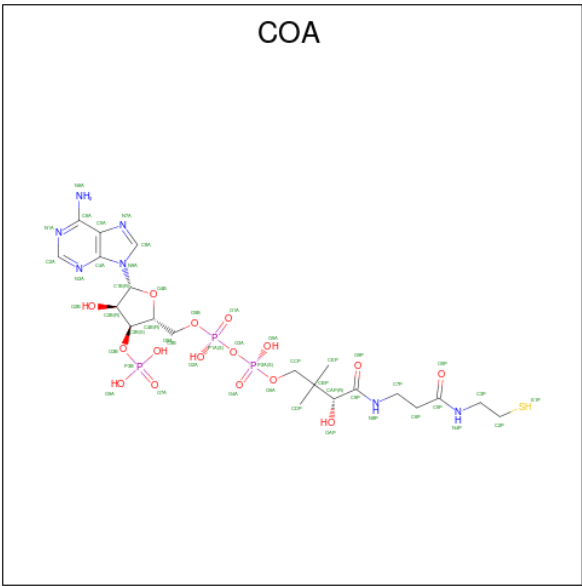
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



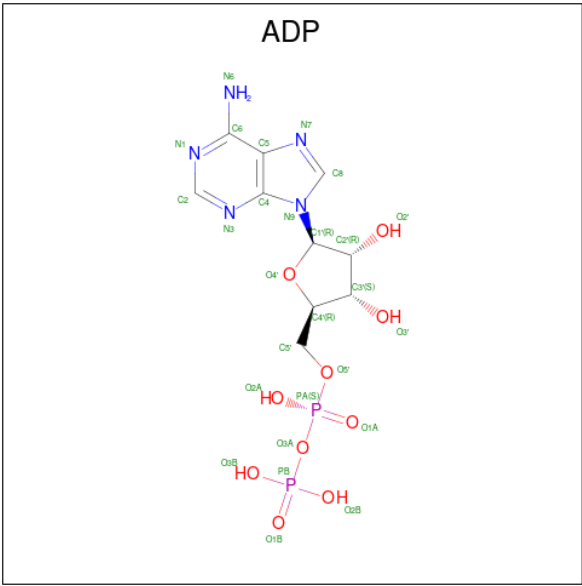
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
5	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 6 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

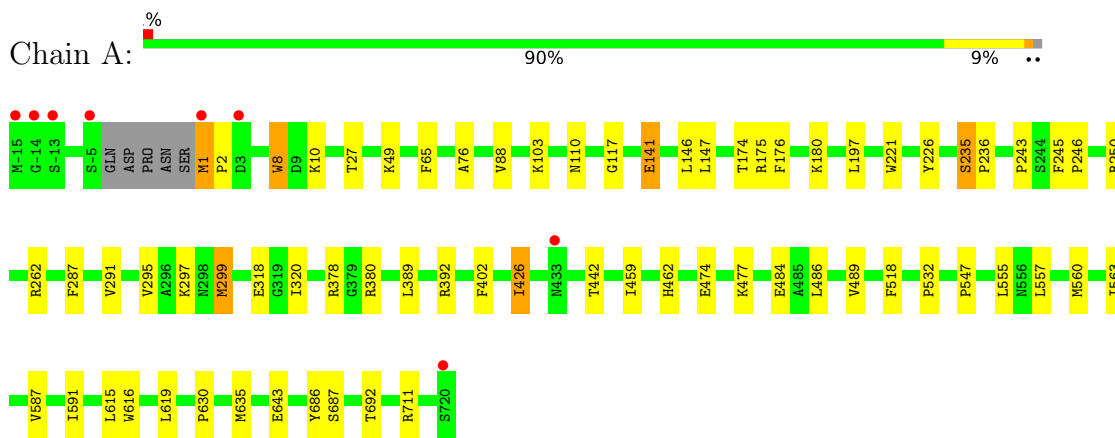
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	235	Total	O	0	0
			235	235		
7	B	254	Total	O	0	0
			254	254		
7	C	168	Total	O	0	0
			168	168		
7	D	122	Total	O	0	0
			122	122		

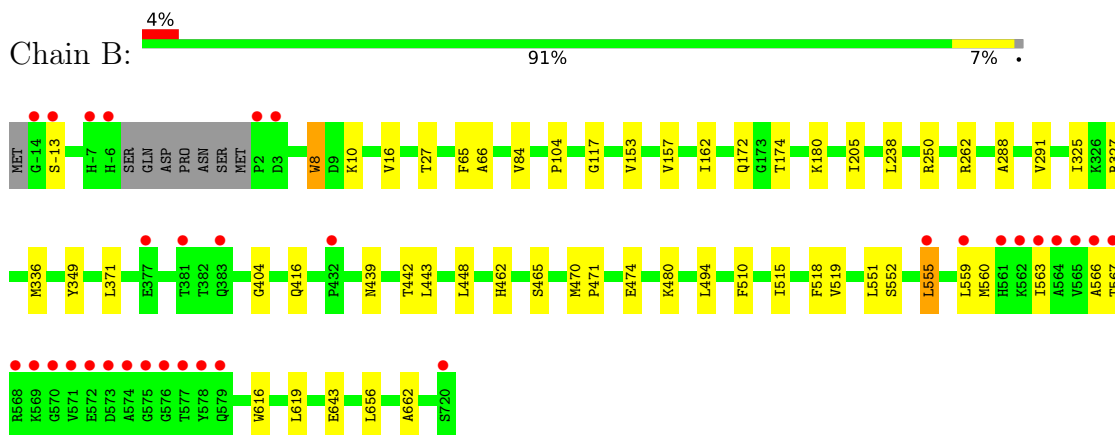
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

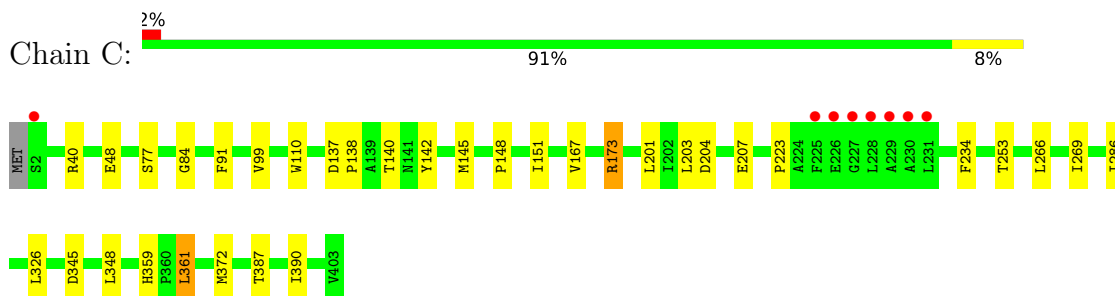
• Molecule 1: FATTY ACID BETA-OXIDATION COMPLEX ALPHA-CHAIN FADB



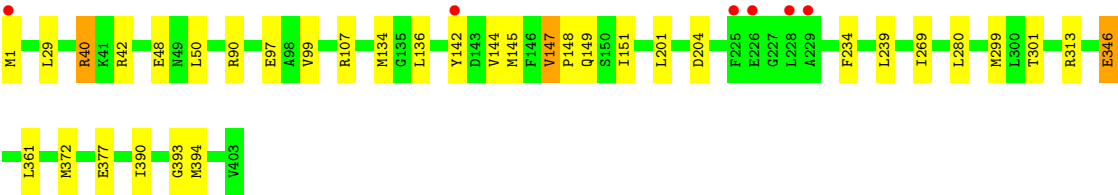
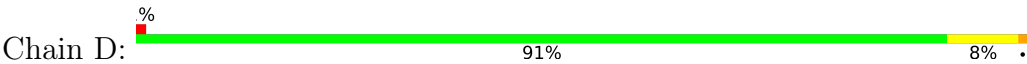
• Molecule 1: FATTY ACID BETA-OXIDATION COMPLEX ALPHA-CHAIN FADB



• Molecule 2: FATTY ACID BETA-OXIDATION COMPLEX BETA-CHAIN FADA



● Molecule 2: FATTY ACID BETA-OXIDATION COMPLEX BETA-CHAIN FADA



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	248.16Å 135.97Å 118.16Å 90.00° 110.43° 90.00°	Depositor
Resolution (Å)	50.00 – 2.63 50.00 – 2.63	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.63) 99.9 (50.00-2.63)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.186 , 0.229 0.187 , 0.230	Depositor DCC
R_{free} test set	5449 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17910	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, SO4, GOL, COA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.55	1/5574 (0.0%)	0.82	2/7547 (0.0%)
1	B	0.55	0/5527	0.83	1/7490 (0.0%)
2	C	0.56	0/3030	0.80	0/4103
2	D	0.56	0/3019	0.80	0/4092
All	All	0.55	1/17150 (0.0%)	0.82	3/23232 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	547	PRO	CA-C	5.75	1.55	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	SER	CA-C-N	6.27	125.76	119.24
1	A	235	SER	C-N-CA	6.27	125.76	119.24
1	B	555	LEU	N-CA-C	-5.10	102.73	110.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5454	0	5477	44	0
1	B	5405	0	5399	43	0
2	C	2976	0	2996	27	0
2	D	2968	0	2975	27	0
3	A	35	0	0	0	0
3	B	45	0	0	0	0
3	C	50	0	0	0	0
3	D	30	0	0	0	0
4	A	12	0	16	2	0
4	B	6	0	8	0	0
5	A	48	0	32	0	0
5	B	48	0	32	0	0
6	A	27	0	12	0	0
6	C	27	0	12	0	0
7	A	235	0	0	3	0
7	B	254	0	0	0	0
7	C	168	0	0	0	0
7	D	122	0	0	0	0
All	All	17910	0	16959	125	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (125) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262[A]:ARG:NE	2:C:142[A]:TYR:CE2	1.98	1.29
1:A:262[B]:ARG:HG3	1:A:262[B]:ARG:HH11	1.06	1.12
1:B:262[A]:ARG:HD3	2:C:142[A]:TYR:CD2	1.86	1.10
1:A:262[B]:ARG:HH11	1:A:262[B]:ARG:CG	1.73	0.99
1:B:262[A]:ARG:CD	2:C:142[A]:TYR:CD2	2.47	0.97
1:A:103:LYS:HE3	7:A:2018:HOH:O	1.66	0.93
1:A:262[B]:ARG:HG3	1:A:262[B]:ARG:NH1	1.80	0.90
1:B:262[A]:ARG:NE	2:C:142[A]:TYR:CD2	2.42	0.87
2:D:90:ARG:HH11	2:D:394:MET:HE1	1.44	0.82
2:C:84:GLY:HA2	2:D:394:MET:HE3	1.62	0.79
1:B:27:THR:HG21	1:B:66:ALA:HB3	1.67	0.75
1:B:84:VAL:HG11	1:B:291:VAL:HG11	1.68	0.74
2:D:40:ARG:HD2	2:D:48:GLU:OE2	1.93	0.69
1:B:416:GLN:HG3	1:B:448:LEU:HD23	1.77	0.67
1:B:563:ILE:HA	1:B:566:ALA:HB3	1.77	0.66
1:B:462:HIS:HB3	1:B:474:GLU:HB3	1.77	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:110:TRP:HD1	2:D:313:ARG:HD3	1.62	0.64
2:C:173:ARG:NH2	2:C:348:LEU:O	2.30	0.64
1:A:287:PHE:CE2	1:A:291:VAL:HG21	2.33	0.64
1:B:262[A]:ARG:HD3	2:C:142[A]:TYR:CG	2.33	0.63
1:B:559:LEU:O	1:B:563:ILE:HG23	1.99	0.63
1:A:711:ARG:HB3	4:A:1728:GOL:H2	1.80	0.63
1:A:262[B]:ARG:CG	1:A:262[B]:ARG:NH1	2.45	0.62
1:A:389:LEU:HA	1:A:392:ARG:NH1	2.15	0.61
2:C:372:MET:HA	2:C:372:MET:HE2	1.83	0.61
1:A:297:LYS:HE3	7:A:2132:HOH:O	2.02	0.59
1:B:27:THR:CG2	1:B:66:ALA:HB3	2.32	0.58
2:C:40:ARG:NH2	2:C:77:SER:O	2.32	0.58
1:A:295:VAL:CG1	1:A:299:MET:HE2	2.33	0.58
1:B:262[A]:ARG:CD	2:C:142[A]:TYR:CE2	2.80	0.58
1:B:8[A]:TRP:CZ3	1:B:10:LYS:HB2	2.40	0.57
1:A:65:PHE:HB3	1:A:117:GLY:HA2	1.87	0.56
1:A:141:GLU:HG3	1:A:147:LEU:C	2.31	0.55
1:B:262[A]:ARG:CZ	2:C:142[A]:TYR:CE2	2.86	0.54
2:D:147:VAL:O	2:D:299:MET:HE1	2.07	0.54
1:A:711:ARG:HH11	4:A:1728:GOL:H12	1.73	0.54
1:B:162:ILE:HD12	1:B:238:LEU:HD21	1.89	0.54
2:C:110:TRP:CH2	2:D:107:ARG:HD2	2.43	0.53
1:A:146:LEU:HD22	1:A:291:VAL:HG22	1.91	0.53
1:A:402:PHE:CD2	1:A:426:ILE:HG12	2.44	0.53
1:B:616:TRP:O	1:B:619:LEU:HB2	2.09	0.53
1:B:65:PHE:HB3	1:B:117:GLY:HA2	1.90	0.53
2:C:84:GLY:CA	2:D:394:MET:HE3	2.37	0.53
2:C:203:LEU:HD11	2:C:207:GLU:HG3	1.91	0.52
1:A:295:VAL:HG12	1:A:299:MET:HE2	1.92	0.52
2:D:201:LEU:HD11	2:D:204:ASP:HB3	1.92	0.52
1:B:552:SER:O	1:B:555:LEU:O	2.28	0.51
2:D:90:ARG:HD3	2:D:394:MET:HE2	1.91	0.51
2:C:201:LEU:HD11	2:C:204:ASP:HB3	1.93	0.51
1:B:515:ILE:HD11	1:B:551:LEU:HD21	1.94	0.50
1:B:336:MET:SD	1:B:465:SER:HB3	2.52	0.50
2:D:134:MET:HB2	2:D:144:VAL:HG21	1.94	0.50
1:A:518:PHE:HB2	1:A:643:GLU:CD	2.37	0.50
2:D:372:MET:HA	2:D:372:MET:HE2	1.94	0.49
1:A:243:PRO:HG3	2:D:136:LEU:HD23	1.94	0.49
1:A:686:TYR:O	1:A:692:THR:HA	2.12	0.49
1:A:8[A]:TRP:CZ3	1:A:10:LYS:HB2	2.47	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:462:HIS:HB3	1:A:474:GLU:HB3	1.94	0.49
1:B:84:VAL:HG11	1:B:291:VAL:CG1	2.39	0.49
1:B:262[A]:ARG:NE	2:C:142[A]:TYR:HE2	1.92	0.48
1:A:8[A]:TRP:HZ3	1:A:10:LYS:HB2	1.78	0.48
2:D:90:ARG:HH11	2:D:394:MET:CE	2.20	0.48
1:A:65:PHE:HB3	1:A:117:GLY:CA	2.43	0.48
1:B:325:ILE:HB	1:B:349:TYR:CE1	2.49	0.47
1:B:518:PHE:HB2	1:B:643:GLU:CD	2.39	0.47
2:D:90:ARG:HH21	2:D:97:GLU:CD	2.21	0.47
1:B:327:ARG:HB3	1:B:404:GLY:O	2.15	0.47
2:D:148:PRO:HD3	2:D:234:PHE:CD1	2.50	0.47
2:C:140:THR:HG21	2:D:29:LEU:HD23	1.96	0.46
1:A:555:LEU:HD12	1:A:560:MET:HE2	1.96	0.46
1:A:442:THR:HG21	1:A:563:ILE:HG12	1.98	0.46
1:A:245:PHE:HB2	1:A:246:PRO:HD3	1.98	0.46
1:B:8[A]:TRP:HZ3	1:B:10:LYS:HB2	1.79	0.46
1:B:250:ARG:NH1	2:C:145:MET:HG2	2.31	0.46
1:A:221:TRP:HA	1:A:226:TYR:CD1	2.50	0.46
1:B:27:THR:HG21	1:B:66:ALA:CB	2.42	0.45
2:C:223:PRO:HA	2:C:253:THR:HG22	1.98	0.45
1:B:442:THR:HG21	1:B:563:ILE:CG2	2.46	0.45
1:A:76:ALA:O	1:A:297:LYS:NZ	2.43	0.45
1:B:656:LEU:HD13	1:B:662:ALA:HB2	1.98	0.45
2:D:390:ILE:HB	2:D:394:MET:HB2	1.97	0.45
1:A:630:PRO:HG2	1:A:635:MET:HE2	1.99	0.45
1:A:459:ILE:HG21	1:A:489:VAL:HG21	1.99	0.44
1:A:587:VAL:O	1:A:591:ILE:HG12	2.17	0.44
1:A:378:ARG:HH21	1:A:380:ARG:NH2	2.15	0.44
7:A:2110:HOH:O	2:D:142[B]:TYR:CZ	2.68	0.44
2:C:40:ARG:HD3	2:C:48:GLU:OE2	2.18	0.44
2:D:299:MET:O	2:D:393:GLY:HA2	2.17	0.44
1:A:616:TRP:O	1:A:619:LEU:HB2	2.17	0.44
1:A:250:ARG:HD2	2:D:142[B]:TYR:CE2	2.53	0.44
1:B:510:PHE:CD1	1:B:656:LEU:HD11	2.53	0.44
2:C:91:PHE:HB2	2:C:390:ILE:HG23	2.00	0.44
2:D:301:THR:O	2:D:301:THR:HG22	2.18	0.44
1:A:378:ARG:HH21	1:A:380:ARG:HH22	1.65	0.43
2:D:1:MET:HA	2:D:107:ARG:HH21	1.82	0.43
1:A:235:SER:HA	1:A:236:PRO:HD3	1.90	0.43
1:B:65:PHE:HB3	1:B:117:GLY:CA	2.49	0.43
1:B:560:MET:HA	1:B:563:ILE:HG12	1.99	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:99:VAL:HG13	2:D:269:ILE:HD11	1.99	0.43
1:A:287:PHE:O	1:A:291:VAL:HG23	2.19	0.43
1:A:616:TRP:HE3	1:A:619:LEU:HD13	1.83	0.43
1:B:470:MET:HA	1:B:471:PRO:HD3	1.79	0.43
2:C:326:LEU:HD13	2:C:387:THR:HG23	2.01	0.43
2:C:99:VAL:HG13	2:C:269:ILE:HD11	2.01	0.42
2:C:359:HIS:CE1	2:C:361:LEU:HA	2.54	0.42
1:B:325:ILE:HB	1:B:349:TYR:HE1	1.83	0.42
1:A:532:PRO:HB2	1:A:615:LEU:HD13	2.02	0.42
1:A:1:MET:HA	1:A:2:PRO:HD3	1.94	0.42
2:D:1:MET:HA	2:D:107:ARG:NH2	2.34	0.42
2:C:148:PRO:HD3	2:C:234:PHE:CD1	2.55	0.41
1:A:320:ILE:HG21	1:A:484:GLU:HA	2.01	0.41
1:A:176:PHE:CD1	1:A:180:LYS:HD2	2.56	0.41
1:B:439:ASN:CG	1:B:439:ASN:O	2.63	0.41
1:A:477:LYS:HB2	1:A:486:LEU:HD21	2.02	0.41
1:B:104:PRO:HG2	1:B:205:ILE:HG23	2.02	0.41
2:D:90:ARG:HD3	2:D:394:MET:CE	2.50	0.41
1:B:563:ILE:HA	1:B:566:ALA:CB	2.49	0.41
2:D:346:GLU:CD	2:D:346:GLU:H	2.29	0.41
1:A:110:ASN:HA	1:A:197:LEU:HD11	2.03	0.41
2:D:149:GLN:HA	2:D:299:MET:HE3	2.02	0.40
1:B:153:VAL:O	1:B:157:VAL:HG23	2.21	0.40
1:B:519:VAL:HG21	1:B:560:MET:HE2	2.04	0.40
2:C:137:ASP:HA	2:C:138:PRO:HD3	1.95	0.40
1:B:288:ALA:HA	1:B:291:VAL:HG12	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	732/736 (100%)	708 (97%)	23 (3%)	1 (0%)	48	64
1	B	730/736 (99%)	710 (97%)	20 (3%)	0	100	100
2	C	403/403 (100%)	390 (97%)	12 (3%)	1 (0%)	43	58
2	D	403/403 (100%)	391 (97%)	11 (3%)	1 (0%)	43	58
All	All	2268/2278 (100%)	2199 (97%)	66 (3%)	3 (0%)	48	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	361	LEU
2	D	361	LEU
1	A	318	GLU

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	560/566 (99%)	547 (98%)	13 (2%)	44	65
1	B	553/566 (98%)	541 (98%)	12 (2%)	45	66
2	C	309/310 (100%)	303 (98%)	6 (2%)	50	70
2	D	307/310 (99%)	297 (97%)	10 (3%)	33	53
All	All	1729/1752 (99%)	1688 (98%)	41 (2%)	44	64

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	8[A]	TRP
1	A	8[B]	TRP
1	A	27	THR
1	A	49	LYS
1	A	88	VAL
1	A	141	GLU
1	A	174	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	175	ARG
1	A	299	MET
1	A	426	ILE
1	A	557	LEU
1	A	687	SER
1	B	-13	SER
1	B	8[A]	TRP
1	B	8[B]	TRP
1	B	16	VAL
1	B	172	GLN
1	B	174	THR
1	B	180	LYS
1	B	371	LEU
1	B	443	LEU
1	B	480	LYS
1	B	494	LEU
1	B	567	THR
2	C	151	ILE
2	C	167	VAL
2	C	173	ARG
2	C	266	LEU
2	C	286	ILE
2	C	345	ASP
2	D	40	ARG
2	D	42	ARG
2	D	50	LEU
2	D	145	MET
2	D	147	VAL
2	D	151	ILE
2	D	239	LEU
2	D	280	LEU
2	D	346	GLU
2	D	377	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	294	GLN
1	A	416	GLN
1	A	561	HIS
1	A	681	GLN
1	B	4	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	163	GLN
1	B	416	GLN
1	B	556	ASN
1	B	633	GLN
2	C	49	ASN
2	D	338	GLN
2	D	380	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

39 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	A	1723	-	4,4,4	0.44	0	6,6,6	0.12	0
3	SO4	B	1729	-	4,4,4	0.44	0	6,6,6	0.16	0
3	SO4	C	1411	-	4,4,4	0.46	0	6,6,6	0.09	0
3	SO4	A	1727	-	4,4,4	0.42	0	6,6,6	0.14	0
3	SO4	C	1407	-	4,4,4	0.45	0	6,6,6	0.10	0
3	SO4	D	1408	-	4,4,4	0.44	0	6,6,6	0.12	0
3	SO4	A	1724	-	4,4,4	0.45	0	6,6,6	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	C	1406	-	4,4,4	0.42	0	6,6,6	0.10	0
3	SO4	C	1405	-	4,4,4	0.44	0	6,6,6	0.06	0
3	SO4	D	1407	-	4,4,4	0.45	0	6,6,6	0.11	0
4	GOL	A	1728	-	5,5,5	0.35	0	5,5,5	0.23	0
4	GOL	B	1730	-	5,5,5	0.36	0	5,5,5	0.38	0
5	COA	A	1730	-	47,50,50	1.22	6 (12%)	69,75,75	1.49	11 (15%)
3	SO4	B	1726	-	4,4,4	0.44	0	6,6,6	0.14	0
3	SO4	B	1728	-	4,4,4	0.45	0	6,6,6	0.11	0
3	SO4	D	1409	-	4,4,4	0.44	0	6,6,6	0.09	0
3	SO4	B	1724	-	4,4,4	0.42	0	6,6,6	0.16	0
6	ADP	C	1414	-	28,29,29	1.59	5 (17%)	43,45,45	1.86	8 (18%)
3	SO4	A	1726	-	4,4,4	0.44	0	6,6,6	0.15	0
4	GOL	A	1729	-	5,5,5	0.36	0	5,5,5	0.31	0
3	SO4	D	1406	-	4,4,4	0.44	0	6,6,6	0.14	0
3	SO4	B	1721	-	4,4,4	0.43	0	6,6,6	0.15	0
3	SO4	C	1404	-	4,4,4	0.40	0	6,6,6	0.26	0
3	SO4	C	1408	-	4,4,4	0.43	0	6,6,6	0.14	0
3	SO4	A	1721	-	4,4,4	0.45	0	6,6,6	0.19	0
3	SO4	C	1410	-	4,4,4	0.41	0	6,6,6	0.20	0
3	SO4	C	1412	-	4,4,4	0.45	0	6,6,6	0.10	0
3	SO4	B	1723	-	4,4,4	0.44	0	6,6,6	0.10	0
3	SO4	A	1725	-	4,4,4	0.44	0	6,6,6	0.15	0
3	SO4	D	1405	-	4,4,4	0.44	0	6,6,6	0.07	0
3	SO4	C	1413	-	4,4,4	0.46	0	6,6,6	0.13	0
3	SO4	D	1404	-	4,4,4	0.44	0	6,6,6	0.10	0
3	SO4	B	1722	-	4,4,4	0.46	0	6,6,6	0.07	0
3	SO4	B	1725	-	4,4,4	0.40	0	6,6,6	0.25	0
5	COA	B	1731	-	47,50,50	1.25	6 (12%)	69,75,75	1.57	12 (17%)
3	SO4	A	1722	-	4,4,4	0.46	0	6,6,6	0.17	0
3	SO4	C	1409	-	4,4,4	0.44	0	6,6,6	0.14	0
6	ADP	A	1731	-	28,29,29	1.46	3 (10%)	43,45,45	1.92	10 (23%)
3	SO4	B	1727	-	4,4,4	0.45	0	6,6,6	0.07	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	COA	B	1731	-	-	12/48/64/64	0/3/3/3
4	GOL	A	1729	-	-	4/4/4/4	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	ADP	C	1414	-	-	2/16/32/32	0/3/3/3
6	ADP	A	1731	-	-	5/16/32/32	0/3/3/3
4	GOL	B	1730	-	-	2/4/4/4	-
4	GOL	A	1728	-	-	2/4/4/4	-
5	COA	A	1730	-	-	5/48/64/64	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	1414	ADP	C5-C4	5.24	1.48	1.39
5	B	1731	COA	C5A-C4A	4.97	1.48	1.39
5	A	1730	COA	C5A-C4A	4.93	1.47	1.39
6	A	1731	ADP	C5-C4	4.91	1.47	1.39
6	C	1414	ADP	PA-O3A	3.24	1.63	1.59
6	A	1731	ADP	C5-C6	3.08	1.49	1.41
6	C	1414	ADP	C5-C6	2.98	1.49	1.41
5	B	1731	COA	C5A-C6A	2.91	1.49	1.41
5	A	1730	COA	C5A-C6A	2.90	1.49	1.41
6	C	1414	ADP	C8-N7	2.51	1.36	1.31
5	B	1731	COA	P1A-O3A	2.50	1.62	1.59
5	B	1731	COA	C8A-N7A	2.41	1.36	1.31
6	A	1731	ADP	C8-N7	2.38	1.36	1.31
5	B	1731	COA	P2A-O3A	2.32	1.62	1.59
5	A	1730	COA	C8A-N7A	2.30	1.36	1.31
5	A	1730	COA	P1A-O3A	2.14	1.61	1.59
5	A	1730	COA	P2A-O3A	2.10	1.61	1.59
5	B	1731	COA	C5A-N7A	-2.10	1.35	1.39
6	C	1414	ADP	C5-N7	-2.06	1.35	1.39
5	A	1730	COA	C5A-N7A	-2.03	1.35	1.39

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1414	ADP	C5-C4-N3	-6.21	118.17	126.72
6	A	1731	ADP	C5-C4-N3	-5.77	118.78	126.72
5	B	1731	COA	C5A-C4A-N3A	-5.61	118.99	126.72
5	A	1730	COA	C5A-C4A-N3A	-5.25	119.48	126.72
6	C	1414	ADP	N3-C4-N9	4.90	135.50	127.17
6	A	1731	ADP	N3-C4-N9	4.62	135.02	127.17
5	B	1731	COA	N3A-C4A-N9A	4.57	134.94	127.17
5	A	1730	COA	N3A-C4A-N9A	4.46	134.76	127.17

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1731	ADP	C2-N3-C4	4.01	121.64	111.83
6	C	1414	ADP	C2-N3-C4	3.91	121.37	111.83
5	B	1731	COA	C2A-N3A-C4A	3.89	121.33	111.83
5	B	1731	COA	N3A-C2A-N1A	-3.81	122.82	128.58
6	A	1731	ADP	N3-C2-N1	-3.77	122.87	128.58
5	A	1730	COA	C2A-N3A-C4A	3.65	120.73	111.83
5	A	1730	COA	N3A-C2A-N1A	-3.62	123.10	128.58
6	A	1731	ADP	C4-C5-N7	-3.56	106.51	110.58
5	B	1731	COA	C4A-C5A-N7A	-3.41	106.68	110.58
6	C	1414	ADP	C4-C5-N7	-3.40	106.70	110.58
6	C	1414	ADP	N3-C2-N1	-3.39	123.45	128.58
5	A	1730	COA	C4A-C5A-N7A	-3.21	106.91	110.58
5	A	1730	COA	C4A-N9A-C8A	3.14	109.03	105.74
5	B	1731	COA	O4B-C1B-N9A	3.10	114.05	108.09
6	A	1731	ADP	O4'-C1'-N9	2.98	113.82	108.09
5	B	1731	COA	C4A-N9A-C8A	2.84	108.72	105.74
6	A	1731	ADP	C4-N9-C8	2.76	108.63	105.74
5	A	1730	COA	O4B-C1B-N9A	2.74	113.36	108.09
6	A	1731	ADP	C5-N7-C8	2.67	107.65	103.45
5	B	1731	COA	CEP-CBP-CAP	2.66	113.30	108.77
5	B	1731	COA	C5A-N7A-C8A	2.56	107.47	103.45
6	C	1414	ADP	C3'-C2'-C1'	2.54	106.27	101.46
6	A	1731	ADP	C6-C5-N7	2.52	136.95	132.09
6	C	1414	ADP	C5-N7-C8	2.47	107.34	103.45
5	A	1730	COA	C5A-N7A-C8A	2.45	107.30	103.45
5	B	1731	COA	C6A-C5A-N7A	2.39	136.69	132.09
5	A	1730	COA	C6A-C5A-N7A	2.29	136.51	132.09
5	A	1730	COA	C2A-N1A-C6A	2.18	122.31	118.73
5	B	1731	COA	C2A-N1A-C6A	2.14	122.25	118.73
5	A	1730	COA	N9A-C8A-N7A	-2.07	111.00	113.94
6	A	1731	ADP	N9-C8-N7	-2.06	111.01	113.94
6	C	1414	ADP	C4-N9-C8	2.05	107.89	105.74
5	B	1731	COA	N9A-C8A-N7A	-2.03	111.05	113.94

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1729	GOL	O1-C1-C2-C3
4	A	1729	GOL	C1-C2-C3-O3
5	A	1730	COA	C6P-C5P-N4P-C3P
5	B	1731	COA	CCP-O6A-P2A-O3A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
5	B	1731	COA	CCP-O6A-P2A-O4A
5	B	1731	COA	CCP-O6A-P2A-O5A
5	B	1731	COA	C9P-CAP-CBP-CCP
5	B	1731	COA	OAP-CAP-CBP-CDP
5	B	1731	COA	C9P-CAP-CBP-CDP
5	B	1731	COA	C9P-CAP-CBP-CEP
5	B	1731	COA	CAP-C9P-N8P-C7P
6	A	1731	ADP	C5'-O5'-PA-O3A
6	A	1731	ADP	C4'-C5'-O5'-PA
5	A	1730	COA	O5P-C5P-N4P-C3P
5	B	1731	COA	O9P-C9P-N8P-C7P
6	A	1731	ADP	C3'-C4'-C5'-O5'
4	B	1730	GOL	C1-C2-C3-O3
4	A	1729	GOL	O1-C1-C2-O2
4	A	1729	GOL	O2-C2-C3-O3
6	A	1731	ADP	O4'-C4'-C5'-O5'
6	C	1414	ADP	O4'-C4'-C5'-O5'
5	B	1731	COA	OAP-CAP-CBP-CEP
4	B	1730	GOL	O2-C2-C3-O3
6	C	1414	ADP	C3'-C4'-C5'-O5'
5	B	1731	COA	P1A-O3A-P2A-O6A
5	B	1731	COA	OAP-CAP-CBP-CCP
6	A	1731	ADP	C5'-O5'-PA-O1A
4	A	1728	GOL	C1-C2-C3-O3
5	A	1730	COA	C3B-O3B-P3B-O7A
4	A	1728	GOL	O2-C2-C3-O3
5	A	1730	COA	C3B-O3B-P3B-O8A
5	A	1730	COA	C3B-O3B-P3B-O9A

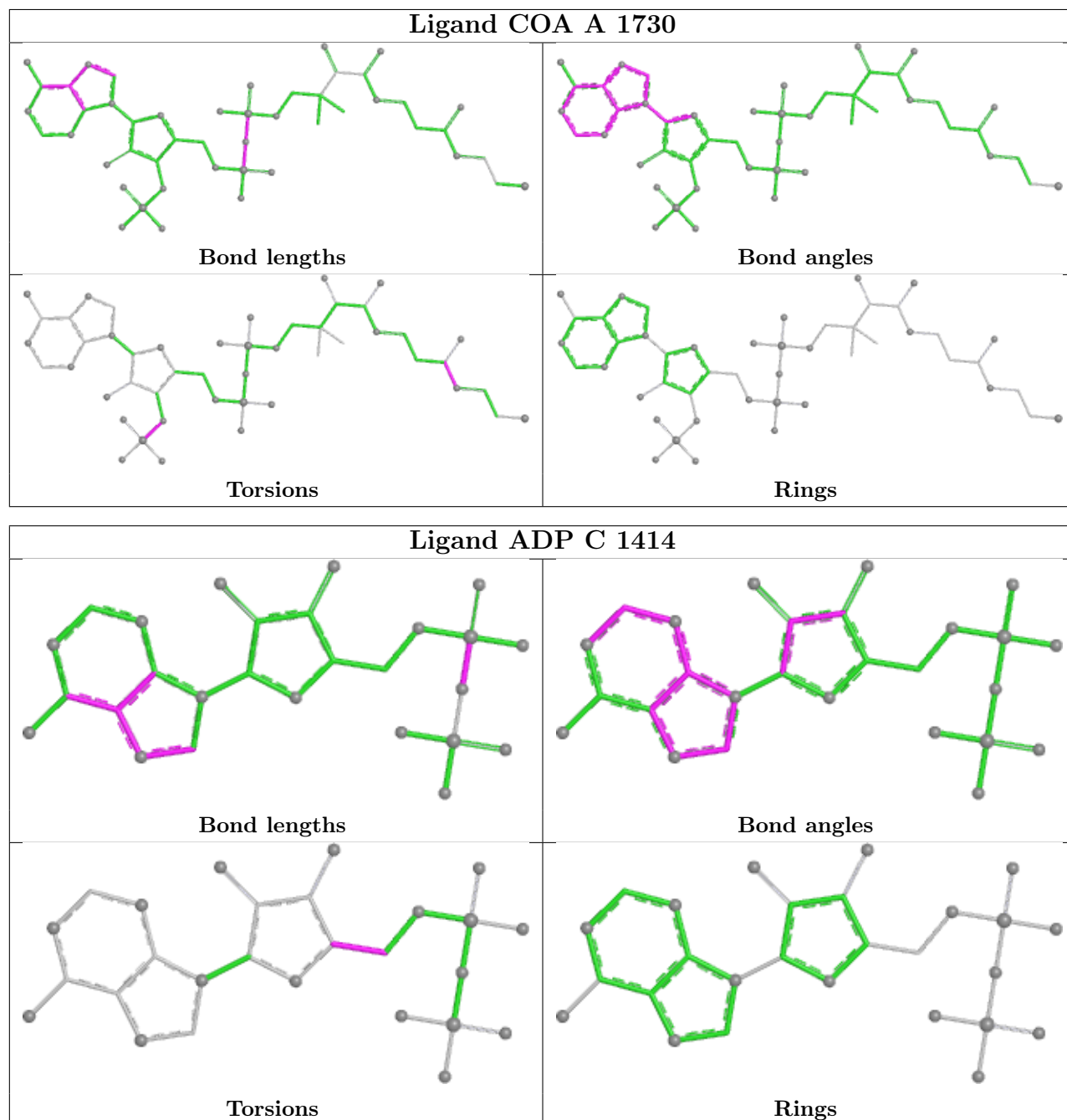
There are no ring outliers.

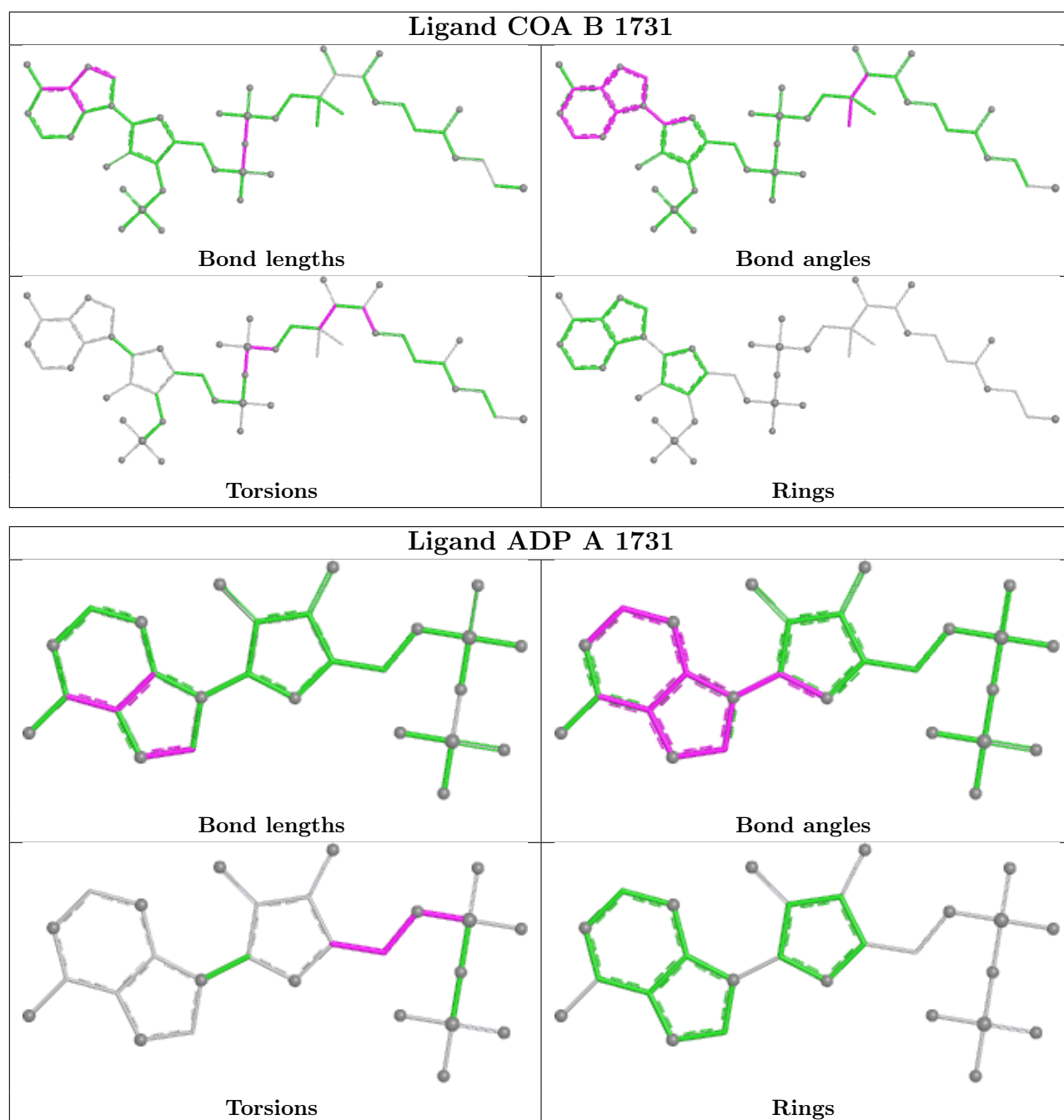
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1728	GOL	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	731/736 (99%)	-0.33	8 (1%) 78 75	15, 40, 63, 90	5 (0%)
1	B	728/736 (98%)	-0.15	32 (4%) 39 32	15, 38, 70, 115	6 (0%)
2	C	402/403 (99%)	-0.55	8 (1%) 65 61	17, 29, 47, 102	3 (0%)
2	D	403/403 (100%)	-0.48	6 (1%) 72 69	16, 34, 51, 84	2 (0%)
All	All	2264/2278 (99%)	-0.34	54 (2%) 59 55	15, 36, 64, 115	16 (0%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-14	GLY	7.6
2	C	229	ALA	7.5
1	B	566	ALA	6.1
2	C	230	ALA	5.6
2	C	225	PHE	5.5
1	B	565	VAL	5.4
1	A	720	SER	5.3
2	C	226	GLU	5.2
1	B	567	THR	4.8
2	C	228	LEU	4.7
1	B	570	GLY	4.6
1	B	574	ALA	4.5
2	C	227	GLY	4.5
1	A	1	MET	4.5
1	B	573	ASP	4.4
1	A	-15	MET	4.3
1	B	563	ILE	4.2
1	B	-6	HIS	3.9
2	D	1	MET	3.8
1	A	-14	GLY	3.7
1	B	571	VAL	3.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	-5	SER	3.7
1	B	2	PRO	3.5
2	D	226	GLU	3.5
1	B	578	TYR	3.4
2	D	229	ALA	3.4
1	B	569	LYS	3.3
1	A	-13	SER	3.3
2	C	2	SER	3.1
2	D	225	PHE	3.1
1	B	564	ALA	2.9
1	B	-13	SER	2.9
1	B	572	GLU	2.8
1	B	559	LEU	2.8
1	B	576	GLY	2.8
2	D	228	LEU	2.6
1	B	575	GLY	2.5
1	B	577	THR	2.5
1	B	-7	HIS	2.5
2	C	231	LEU	2.5
1	B	432	PRO	2.5
1	B	720	SER	2.4
1	B	568	ARG	2.3
1	B	561	HIS	2.2
2	D	142[A]	TYR	2.2
1	A	3	ASP	2.2
1	B	383	GLN	2.2
1	B	579	GLN	2.2
1	B	3	ASP	2.2
1	A	433	ASN	2.2
1	B	377	GLU	2.1
1	B	555	LEU	2.1
1	B	562	LYS	2.1
1	B	381	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	D	1409	5/5	0.62	0.22	93,94,96,97	5
6	ADP	A	1731	27/27	0.64	0.20	92,105,119,121	0
6	ADP	C	1414	27/27	0.72	0.19	67,87,109,110	0
3	SO4	A	1726	5/5	0.75	0.17	48,48,49,50	5
3	SO4	A	1725	5/5	0.76	0.15	39,40,40,41	5
3	SO4	B	1727	5/5	0.79	0.26	30,30,31,31	5
3	SO4	B	1723	5/5	0.80	0.13	62,63,64,64	5
3	SO4	A	1724	5/5	0.80	0.14	66,67,69,69	5
3	SO4	B	1729	5/5	0.80	0.17	85,86,87,88	0
3	SO4	B	1724	5/5	0.81	0.14	43,44,44,44	5
3	SO4	C	1412	5/5	0.83	0.21	47,48,49,49	5
3	SO4	A	1727	5/5	0.85	0.28	23,23,23,23	5
4	GOL	B	1730	6/6	0.87	0.16	53,54,54,54	0
3	SO4	B	1725	5/5	0.87	0.17	57,57,58,58	5
3	SO4	C	1411	5/5	0.87	0.11	47,48,48,49	5
5	COA	B	1731	48/48	0.88	0.14	66,76,82,83	0
3	SO4	C	1410	5/5	0.89	0.11	44,46,46,47	5
3	SO4	C	1409	5/5	0.90	0.13	49,49,51,53	5
4	GOL	A	1729	6/6	0.90	0.15	48,50,50,51	0
4	GOL	A	1728	6/6	0.91	0.10	54,55,56,57	0
3	SO4	C	1407	5/5	0.91	0.20	72,74,76,77	0
3	SO4	B	1728	5/5	0.91	0.19	62,63,63,64	0
3	SO4	B	1726	5/5	0.92	0.11	46,46,48,48	5
3	SO4	D	1406	5/5	0.92	0.16	61,62,64,65	0
5	COA	A	1730	48/48	0.92	0.10	49,59,65,67	0
3	SO4	C	1406	5/5	0.93	0.17	65,67,68,69	0
3	SO4	D	1408	5/5	0.93	0.13	64,64,66,66	0
3	SO4	A	1723	5/5	0.93	0.12	67,68,69,69	0
3	SO4	C	1413	5/5	0.93	0.16	58,61,64,64	0
3	SO4	D	1404	5/5	0.93	0.09	70,72,73,73	0
3	SO4	C	1408	5/5	0.94	0.10	51,52,53,53	5
3	SO4	D	1405	5/5	0.95	0.17	59,60,61,61	0
3	SO4	A	1722	5/5	0.96	0.13	47,49,51,51	0
3	SO4	B	1721	5/5	0.96	0.10	50,51,51,52	0
3	SO4	B	1722	5/5	0.96	0.14	55,55,56,56	0
3	SO4	C	1405	5/5	0.97	0.15	48,49,50,51	0
3	SO4	C	1404	5/5	0.97	0.08	37,40,41,41	0

Continued on next page...

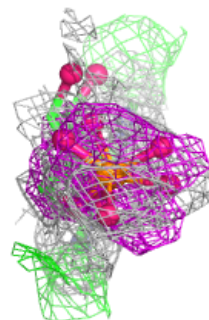
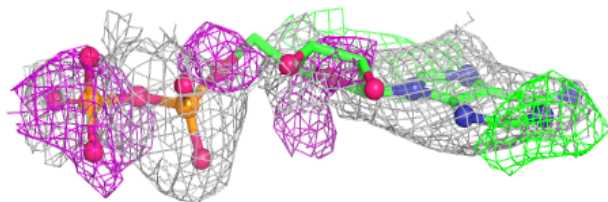
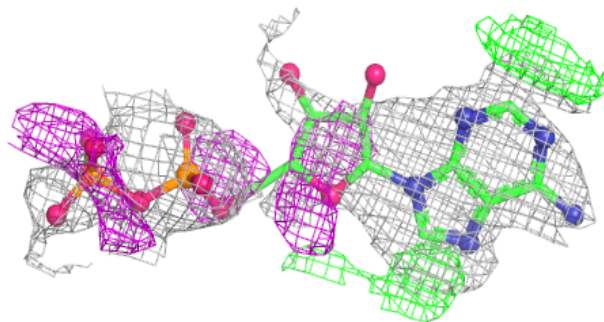
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	D	1407	5/5	0.97	0.11	59,60,60,60	0
3	SO4	A	1721	5/5	0.98	0.09	46,47,47,48	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

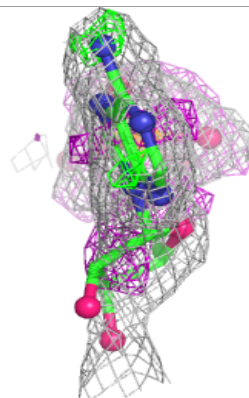
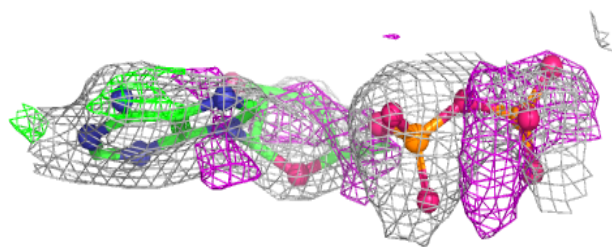
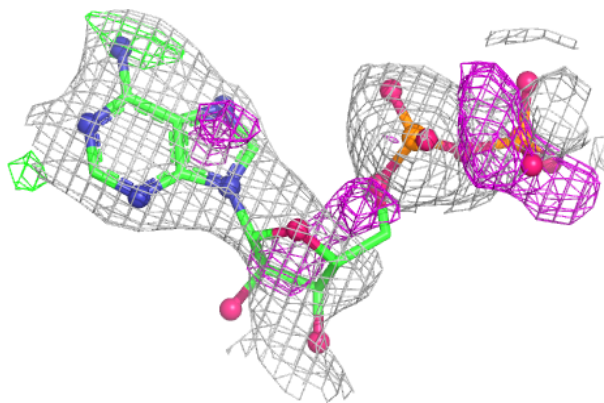
Electron density around ADP A 1731:

2mF_o-DF_c (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
 and green (positive)



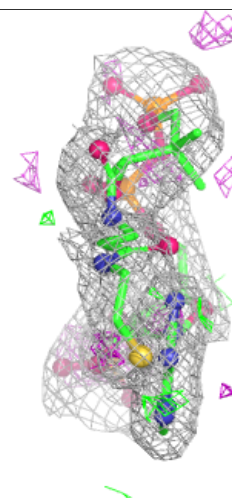
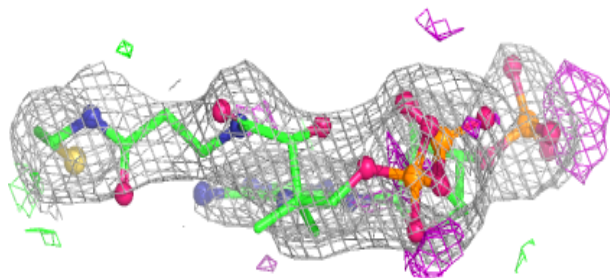
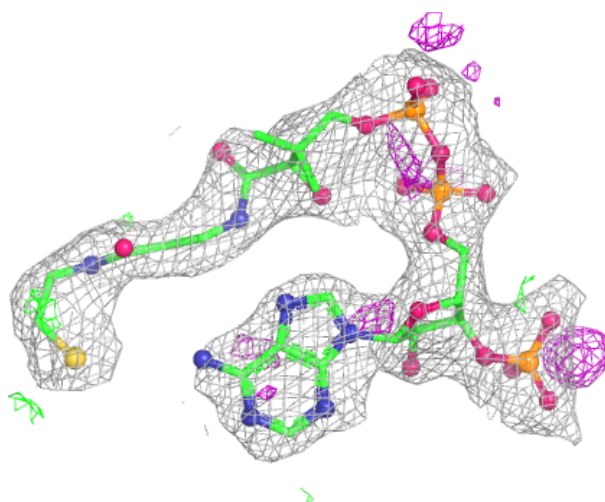
Electron density around ADP C 1414:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



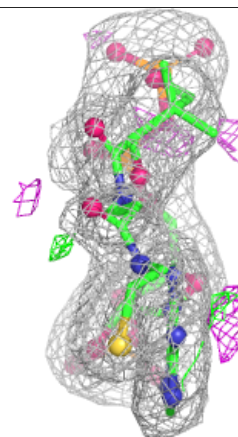
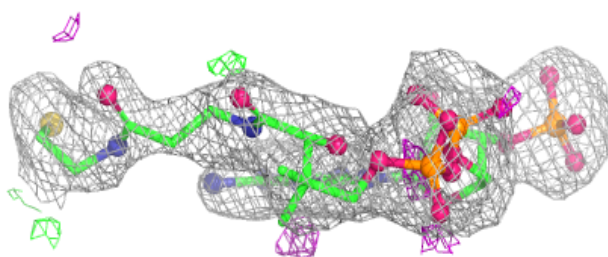
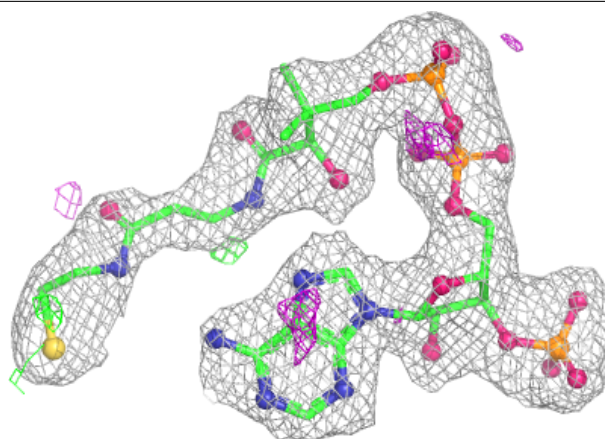
Electron density around COA B 1731:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around COA A 1730:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.